

Information content of the non-linear power spectrum: the effect of beat-coupling to large scales

Christopher D. Rimes^{1*} and Andrew J. S. Hamilton^{1,2*}

¹*JILA, University of Colorado, 440 UCB, Boulder, CO 80309-0440, U.S.A.*

²*Department of Astrophysical and Planetary Sciences, University of Colorado, 391 UCB, Boulder, CO 80309-0391, U.S.A.*

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ABSTRACT

We measure the covariance of the non-linear matter power spectrum from N -body simulations using two methods. In the first case, the covariance of power is estimated from the scatter over many random realizations of the density field. In the second, we use a novel technique to measure the covariance matrix from each simulation individually by re-weighting the density field with a carefully chosen set of functions. The two methods agree at linear scales, but unexpectedly they disagree substantially at increasingly non-linear scales. Moreover, the covariance of non-linear power measured using the re-weightings method changes with box size. The numerical results are consistent with an explanation given in a companion paper, which argues that the cause of the discrepancy is beat-coupling, in which products of Fourier modes separated by a small wavevector couple by gravitational growth to the large-scale beat mode between them. We calculate the information content of the non-linear power spectrum (about the amplitude of the initial, linear power spectrum) using both methods and confirm the result of a previous paper, that at translinear scales the power spectrum contains little information over and above that in the linear power spectrum, but that there is a marked increase in information at non-linear scales. We suggest that, in real galaxy surveys, the covariance of power at non-linear scales is likely to be dominated by beat-coupling to the largest scales of the survey and that, as a result, only part of the information potentially available at non-linear scales is actually measurable from real galaxy surveys.

Key words: cosmology: theory – large-scale structure of Universe.

1 INTRODUCTION

Recent progress in cosmological parameter estimation has been characterized by rapid convergence of constraints from a wealth of different types of observations – galaxy clustering, the Lyman- α forest power spectrum, galaxy cluster abundances, high-redshift type Ia supernovae, weak gravitational lensing, big-bang nucleosynthesis and many others – toward a single, well-determined ‘concordance’ cosmological model. Of particular importance has been the combination of high precision maps of anisotropies in the cosmic microwave background (CMB) with measurements of galaxy clustering from large redshift surveys, which yield complementary constraints on key cosmological parameters (e.g. Efstathiou et al. 2002; Spergel et al. 2003; Tegmark et al. 2004; Seljak et al. 2005; Sanchez et al. 2005).

Galaxy clustering analyses (e.g. Percival et al. 2001; Tegmark et al. 2004; Cole et al. 2005; Eisenstein et al. 2005)

are generally restricted to scales $\gtrsim 20 h^{-1}$ Mpc, where density fluctuations are still linear, galaxy-matter bias appears to be independent of scale (although it does depend on luminosity and on galaxy type) and the matter power spectrum directly traces the spectrum of density fluctuations at recombination. At smaller scales, where much of the observational data in galaxy surveys lie, the extent to which the linear power spectrum can be recovered from the non-linear power spectrum remains unknown. Non-linear evolution changes the shape of the power spectrum in a non-trivial way, and introduces broad correlations between measurements of power at different wavenumbers (Meiksin & White 1999; Scoccimarro, Zaldarriaga & Hui 1999; Cooray & Hu 2001). The early success of analytic formalisms at producing invertible one-to-one mappings between linear and non-linear spectra (Hamilton et al. 1991; Peacock & Dodds 1994, 1996) suggested that information in the linear power spectrum may be preserved into the non-linear regime, but other work (Meiksin, White & Peacock 1999; Seo & Eisenstein 2005) has shown that non-linear evolution

* E-mail: rimes@colorado.edu, Andrew.Hamilton@colorado.edu

erases features in the power spectrum, perhaps leading to an irreversible loss of information.

In an earlier letter (Rimes & Hamilton 2005, hereafter Paper I), we reported measurements of the amount of information in the non-linear power spectrum about the amplitude of the linear power spectrum for the currently favoured (concordance) cosmological model, from a large ensemble of N -body simulations. We have since discovered a small error in our calculations. In Fig. 7 of the present paper we present a revised version of the results from Paper I. Our conclusions remain unchanged: namely, that there exists little independent information in the translinear regime ($k \simeq 0.2\text{--}0.8\ h\text{Mpc}^{-1}$ at the present day) over and above that in the linear power spectrum, but that in the fully non-linear regime there appears to be a significant amount of information beyond that measurable from the linear power spectrum.

Measuring the information content of the non-linear power spectrum involves measuring the covariance matrix of non-linear power, which, if determined as in Paper I from the scatter over an ensemble, requires performing a large number of N -body simulations. This is costly in terms of computing time, especially if the simulations are of high quality. In an attempt to reduce the computational overhead and to streamline the measurement of information, we devised a new technique, described in detail in a companion paper (Hamilton, Rimes & Scoccimarro 2005, hereafter HRS), for estimating the covariance of power from individual simulations, by re-weighting the density field using a set of carefully chosen windows.

In the present paper, we use the re-weighting technique to measure the amount of information in the non-linear power spectrum, for the same set of simulations used in Paper I. We compare the results to those obtained with the ensemble method and present a number of tests of the method.

Unexpectedly, we find that, far from agreeing, the covariance of power measured by the re-weightings method substantially exceeds that measured by the ensemble method. When we first encountered this discrepancy, it seemed to us that it must be caused by a ‘bug’ in our code, and we performed numerous numerical tests to track it down. Belatedly, we realized that the discrepancy was caused not by a bug but by a real physical process, which we term ‘beat-coupling’. The physical origin of beat-coupling is described by HRS, who illustrate its effects with examples using perturbation theory and the hierarchical model.

This paper is organized as follows. In Section 2 we set out our definition of information, and discuss the decorrelation choices that must be made to allocate information to prescribed wavebands. This section also provides a more detailed exposition of the techniques employed in Paper I. Section 3 describes the numerical simulations used in both this paper and the previous one. Section 4 compares measurements of the covariance of power using both ensemble and re-weightings methods, and describes several tests of the results. In Section 5 we compare the information content of the non-linear power spectrum measured using the two different methods. Our conclusions are summarized in Section 6.

2 INFORMATION

2.1 Fisher information

The Fisher information matrix is defined (e.g. Tegmark, Taylor & Heavens 1997) as

$$F^{\alpha\beta} \equiv - \left\langle \frac{\partial^2 \ln \mathcal{L}}{\partial p_\alpha \partial p_\beta} \right\rangle, \quad (1)$$

where $\mathcal{L}(p_\alpha|\text{data})$ is the likelihood function – the multivariate probability distribution of the model parameters p_α given the available data and a set of model assumptions (the Bayesian prior). Fisher information is additive over independent measurements, clearly a desirable property for information to possess. Its importance in parameter estimation is encapsulated in the Cramér-Rao inequality, which limits the maximum precision with which a single parameter p_α can be measured to

$$\langle \Delta \hat{p}_\alpha^2 \rangle \geq 1/F_{\alpha\alpha}, \quad (2)$$

if the estimator \hat{p}_α is unbiased and if this is the only parameter being estimated from the data. Here and throughout this paper we use hats to distinguish an *estimate* of a quantity from its true value. If estimates of the parameters are Gaussian distributed about their expectation values – a good approximation in the limit of a large amount of data, thanks to the central limit theorem – then their covariance matrix is well-approximated by the inverse of the Fisher matrix:

$$\langle \Delta \hat{p}_\alpha \Delta \hat{p}_\beta \rangle \simeq (F^{-1})_{\alpha\beta}. \quad (3)$$

2.2 Power spectrum

We consider a statistically homogeneous and isotropic density field $\rho(\mathbf{r})$. The power spectrum $P(k)$ of density fluctuations of such a field is defined by

$$\langle \delta_{\mathbf{k}} \delta_{\mathbf{k}'} \rangle = (2\pi)^3 \delta_{3D}(\mathbf{k} + \mathbf{k}') P(k), \quad (4)$$

where $\delta_{\mathbf{k}}$ is the Fourier transform of the overdensity $\delta(\mathbf{r}) \equiv \rho(\mathbf{r})/\bar{\rho} - 1$ and $\delta_{3D}(\mathbf{k})$ is a 3-dimensional Dirac delta function. Statistical isotropy requires that the power spectrum be a function only of the magnitude $k \equiv |\mathbf{k}|$ of the wavevector \mathbf{k} .

For Gaussian fluctuations, each $\delta_{\mathbf{k}}$ has real and imaginary components that are independently Gaussianly distributed with variance $P(k)/2$. Usually, power is estimated by averaging over shells in Fourier space. For Gaussian fluctuations, the expected covariance matrix of shell-averaged estimates of power is diagonal, with variance

$$\langle \Delta \hat{P}(k)^2 \rangle = 2P(k)^2/N_k, \quad (5)$$

where N_k is the number of modes in the shell around k (a finite number in the case of a realistic galaxy survey or a periodic N -body simulation). Here $\delta_{\mathbf{k}}$ and its complex conjugate $\delta_{-\mathbf{k}}$ are counted as contributing two distinct modes, the real and imaginary parts of $\delta_{\mathbf{k}}$.

2.3 Information in the power spectrum

Here, as in Paper I, we measure the Fisher information I in a single parameter: the log of the amplitude A of the initial (post-recombination) matter power spectrum, that is,

$$I \equiv - \left\langle \frac{\partial^2 \ln \mathcal{L}}{\partial \ln A^2} \right\rangle. \quad (6)$$

For Gaussian density fluctuations, the power spectrum completely specifies the statistical properties of the density field, so that the only explicit dependence of the likelihood \mathcal{L} is on the power spectrum. For non-Gaussian fluctuations, the likelihood function may also depend explicitly on other parameters. However, as in Paper I, we consider only the information contained in the power spectrum $P(k)$, in which case the information I defined by equation (6) can be expanded as

$$I = - \left\langle \sum_{k,k'} \frac{\partial \ln P(k)}{\partial \ln A} \frac{\partial^2 \ln \mathcal{L}}{\partial \ln P(k) \partial \ln P(k')} \frac{\partial \ln P(k')}{\partial \ln A} \right\rangle. \quad (7)$$

The middle term on the right-hand side of equation (7) is the Hessian of the vector $\ln P(k)$ of log non-linear powers, the expectation value of which is the Fisher matrix. The power spectrum $P(k)$ is averaged over spherical shells in k -space; a typical shell in the non-linear regime contains several thousand to hundreds of thousands of distinct Fourier modes, so it is reasonable to invoke the central limit theorem to assert that estimates of power will be Gaussianly distributed about their expectation values. This assertion holds even if the density field is itself non-Gaussian. We test this is explicitly in Section 4.2.3. In the Gaussian approximation, the Hessian in equation (7) can be approximated by the inverse of the covariance matrix of estimates of log-power.

The remaining terms in equation (7) are two partial derivatives which describe the sensitivity of the non-linear power to changes in the amplitude A of the initial linear power. In the linear regime these derivatives are identically unity, since $P_L(k) \propto A$; at non-linear scales they are equal to the growth rate of the non-linear power spectrum relative to the linear, which can be conveniently measured from simulations.

The information I has a particularly simple interpretation for Gaussian fluctuations. Following equation (5), it is equal to half the total number N of Gaussian modes:

$$I = N/2. \quad (8)$$

As was found in Paper I and is confirmed in Section 5 of the present paper, the information in the non-linear power spectrum $P(k)$ is significantly less than the information in the linear power spectrum at the same wavenumber. This decrease in information could result from a transfer of information from larger to smaller scales, a diversion of information into other quantities (such as the bispectrum), or an irreversible loss of information. In Paper I we argued that complete loss of information during translinear evolution is inconsistent with our finding that the total amount of information on non-linear scales is increasing with time. The remaining two scenarios could, in principle, be distinguished by measuring the information the bispectrum and higher order statistics but this becomes progressively more difficult as the order increases.

2.4 Decorrelated band powers

The quantity I defined by equation (7) is the total amount of information contained in the non-linear power spectrum

about the parameter $\ln A$. Some of this information is degenerate between measurements of power at different wavenumbers as a result of the broad correlations introduced by non-linear evolution.

Hamilton & Tegmark (2000) showed how to decorrelate a power spectrum by defining a set of windowed band-power estimates. Decorrelation is the process of assigning shared information uniquely to a given wavenumber. Here, we extend their method to the case where we want to decorrelate, not the power spectrum itself, but estimates of some parameter – in this case $\ln A$ – made from the power spectrum.

Following Hamilton & Tegmark (2000), we define our windowed band-power estimates \hat{B}_k by:

$$\frac{\hat{B}_k}{P_k} = \sum_{k'} W_{kk'} \frac{\hat{P}_{k'}}{P_{k'}}, \quad (9)$$

where we use the index notation $P_k = P(k)$ to emphasize the fact that the shell-averaged power spectrum $P(k)$ can be viewed as a discrete vector in Fourier space. The band-power windows $W_{kk'}$ in equation (9) are elements of a decorrelation matrix, each column of the matrix being a discrete window for one band-power B_k . There are many (actually an infinite number) of schemes for decorrelating the power spectrum, corresponding to different ways of sharing out degenerate information between wave bands. The reader is directed to Hamilton & Tegmark (2000) for a discussion of the relative merits of selected decorrelation schemes.

Both sides of equation (9) are scaled by P_k , which is a fiducial power spectrum. This scaling ensures that a given band power B_k is not dominated by leakage from wavenumbers k' where the window is small but $P_{k'}$ is large. The choice $P_k = \langle \hat{P}_k \rangle$ guarantees that the expectation value of the windowed band power estimates at each wavenumber is equal to the original power spectrum, provided that the windows are suitably normalized:

$$\sum_{k'} W_{kk'} = 1. \quad (10)$$

In order that the final estimates of $\ln A$ are uncorrelated, the band-power windows must satisfy

$$\mathbf{W}^\top \mathbf{\Lambda} \mathbf{W} = \mathbf{D} \mathbf{F} \mathbf{D}, \quad (11)$$

where \mathbf{F} is the Fisher matrix of the scaled power spectrum and \mathbf{D} is a diagonal matrix with diagonal elements

$$D_{ii} = \frac{\partial \ln P_{k_i}}{\partial \ln A}. \quad (12)$$

The matrix $\mathbf{D} \mathbf{F} \mathbf{D}$ can be interpreted as the Fisher matrix of estimates of $\ln A$ from the power in different wavebands and it is these estimates (rather than the estimates of power themselves) that we want to decorrelate. We experimented with various decorrelation matrices, eventually opting for the upper triangular matrix \mathbf{U} obtained from a generalized form of the Cholesky decomposition:

$$\mathbf{U}^\top \mathbf{\Lambda} \mathbf{U} = \mathbf{D} \mathbf{F} \mathbf{D}, \quad (13)$$

where $\mathbf{\Lambda}$ is a diagonal matrix – the Fisher matrix of the decorrelated estimates of $\ln A$. Note that this is *not* the same as decorrelating the power spectrum using the Cholesky decomposition of \mathbf{F} and then writing the Fisher matrix of the decorrelated estimates as $\mathbf{D} \mathbf{A} \mathbf{D}$ – as we did in Paper I – because \mathbf{D} does not commute with \mathbf{U} . The two approaches are

only approximately equivalent in the case where the band-power windows are narrow or the elements of \mathbf{D} , given by equation 12, are a slowly varying function of k . The former is a particularly poor approximation, because of the broad correlations present in the non-linear power spectrum. In Section 5, we present a corrected version of the relevant figure (fig. 3) from Paper I. Our conclusions are not altered significantly.

In the central-limit-theorem approximation that estimates of power are Gaussianly distributed, the Fisher matrix \mathbf{F} is approximately equal to the scaled inverse covariance of power:

$$\mathbf{F} \simeq \mathbf{P} \langle \Delta \hat{\mathbf{P}} \Delta \hat{\mathbf{P}}^\top \rangle^{-1} \mathbf{P}^\top. \quad (14)$$

Here, \mathbf{P} is a diagonal matrix whose non-zero elements are equal to the fiducial power spectrum P_k . Mathematically, upper Cholesky decorrelation is equivalent to taking a matrix composed of all the elements of the covariance matrix up to some wavenumber k_{\max} , inverting this matrix, and summing all the elements of the resulting Fisher matrix to arrive at a measure of the accumulated information $I(\leq k_{\max})$ up to that wavenumber.

Upper Cholesky decorrelation yields band-power windows that are highly asymmetric, with the band power at each wavenumber k containing contributions only from power on larger scales. A problem with the other schemes that we tried (including the square root of the scaled Fisher matrix, recommended by Hamilton & Tegmark 2000), is that there is an appreciable covariance between large, linear scales and small, non-linear scales. Applying anything other than upper Cholesky decorrelation assigns some of this covariance to large scales, causing the information at linear scales to depart from the expected Gaussian information, equation (5).

While there is a certain arbitrariness about choosing upper Cholesky decorrelation over other possibilities, the resulting cumulative information $I(\leq k_{\max})$ has the virtue of a simple interpretation: it is the information in the power spectrum $P(k)$ at wavenumbers $k \leq k_{\max}$, uncontaminated by information in power at smaller scales.

Because the Fisher matrix of the uncorrelated band powers $B(k) = B_k$ is by definition diagonal, equation (7) reduces to a sum over a single wavenumber and the cumulative information is:

$$I(\leq k) = - \left\langle \sum_{k=0}^{k_{\max}} \frac{\partial \ln B(k)}{\partial \ln A} \frac{\partial^2 \ln \mathcal{L}}{\partial \ln B(k)^2} \frac{\partial \ln B(k)}{\partial \ln A} \right\rangle. \quad (15)$$

3 SIMULATIONS

The simulations used in this paper are the same as those used in Paper I. The main ensemble comprises 600 gravitational N -body simulations of the concordance cosmological model: 400 simulations with a box size of $256 h^{-1}$ Mpc and a further 200 with a box size of $128 h^{-1}$ Mpc. These simulations were evolved using a particle-mesh (PM) code with 128^3 dark matter particles and a 256^3 force mesh.

An additional 25 simulations, also with a box size of $128 h^{-1}$ Mpc, were run using a parallel version of the adaptive mesh refinement (AMR) code ART

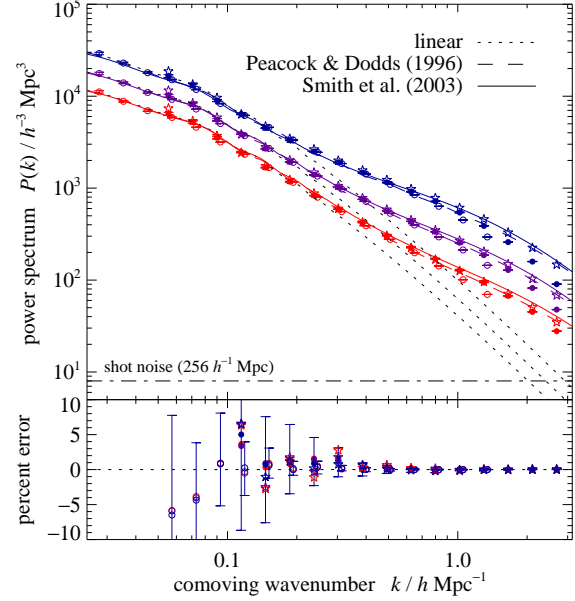


Figure 1. Evolution of the non-linear power spectrum. Top panel: mean power spectrum from the $256 h^{-1}$ Mpc PM simulations (open points, with error bars derived from the scatter between individual simulations); the $128 h^{-1}$ Mpc PM simulations (filled points with error bars); and the $128 h^{-1}$ Mpc ART simulations (stars). Power spectra are shown for three epochs (bottom to top: $a = 0.5, 0.67$ and 1). The linear power spectrum is shown by the dotted curves in each panel. The solid and dashed curves are, respectively, from the fitting formulae of Smith et al. (2003) and Peacock & Dodds (1996). The dot-dashed line marks the level of the shot noise in the $256 h^{-1}$ Mpc simulations; the shot noise in the $128 h^{-1}$ Mpc simulations is a factor of 8 lower. Bottom panel: deviation between the mean power spectra of weighted densities and the power spectrum of the unweighted density. The points are medians from 100 of the PM simulations of each box size and the 25 ART simulations at the same 3 epochs. Error bars mark the upper and lower quartiles of the distribution.

(Kravtsov, Klypin & Khokhlov 1997). Alone, this is an insufficient number to give precise statistics, but together with the much larger ensemble of PM simulations, the ART simulations serve as a useful check of our results on small scales, where the AMR technique is more accurate. For the ART simulations, we used 128^3 particles and a 128^3 root mesh with, at most, three levels of refinement, giving a maximum spatial resolution equivalent to that of a 1024^3 mesh in dense regions. Gaussian initial conditions for the ART simulations were set up using the publicly available GRAFIC package.

The cosmological parameters adopted are those of Tegmark et al. (2004, second-last column of their table 4):

$$(\Omega_M, \Omega_\Lambda, f_b, h, \sigma_8) = (0.29, 0.71, 0.16, 0.71, 0.97) \quad (16)$$

This choice of parameters gives the best fit to the combination of the power spectrum of fluctuations in the cosmic microwave background as measured by the Wilkinson Microwave Anisotropy Probe (WMAP) and galaxy clustering as measured by the Sloan Digital Sky Survey (SDSS), under the important assumptions (among others) of a spatially flat universe ($\Omega_k = 0$) with a cosmological constant ($w = -1$) and a scale-invariant primordial power spectrum

($n_s = 1$). Each simulation was seeded with a different, randomly chosen realization of a Gaussian random field with a power spectrum corresponding to the above cosmological model. The matter transfer function was calculated from the fitting formula of Eisenstein & Hu (1998) for universes with a significant baryon content.

The non-linear power spectra of the above simulations are plotted in Fig. 1 for three epochs: $a = 0.5, 0.67$ and 1.0 , with error bars showing the scatter between individual realizations. Power spectra were calculated by Fourier transforming the weighted density field on a single 256^3 grid¹. The resulting power spectra were corrected for smoothing by dividing by the square of the Fourier transform of the mass assignment window, prior to subtracting the shot noise contribution (Smith et al. 2003). As is to be expected, the power in the PM simulations falls significantly below that measured in the higher resolution ART simulations at small scales, as a result of mesh effects in the PM simulations. The power spectrum from the ART simulations, on the other hand, agrees well with the fits to previous N -body simulations by Peacock & Dodds (1996) and Smith et al. (2003) at the scales of interest. We restrict our analyses to scales for which corrections to power from particle-cell assignment are small, and the shot noise sub-dominant, so uncertainties in both of these corrections should not affect our results.

In what follows, where the results from the $128 h^{-1}$ Mpc PM and ART simulations are consistent, we combine them into a single data set for clarity.

4 CONSTRUCTING THE COVARIANCE MATRIX

Estimating the amount of information in a set of measurements requires knowledge of their covariance matrix. The most direct way to measure the covariance of estimates of the power spectrum is to run an ensemble of N -body simulations, each having a different, random realization of the initial density field (e.g. Meiksin & White 1999; Scoccimarro et al. 1999; Paper I) – a computationally expensive endeavour because many hundreds of realizations are required to yield an accurate estimate of the covariance matrix (Meiksin & White 1999; Paper I).

Alternative approaches to measuring covariances include ‘jackknife’, in which ensembles are formed by removing selected data from the original sample, and ‘bootstrap’, in which ensembles are formed by resampling with replacement (Künsch 1989). These methods work provided that the data being sampled comprise independent random variables drawn from the same distribution. For correlated data,

¹ ‘Chaining’ (Jenkins et al. 1998), a clever way to extend measurements of the power spectrum to smaller scales by superposing the eight octants of a periodic cube on to a single octant, unfortunately cannot be applied to the measurement of covariance, because it involves a reduction in the number of Fourier modes. For Gaussian fluctuations, each mode is independent, so each folding reduces the number of modes by a constant factor 8. At non-linear scales, however, adjacent modes are highly correlated, so subsampling them by a factor of 8 does not reduce the effective number of modes correspondingly.

Künsch (1989) suggested an extension to the bootstrap approach, in which the data are first split into blocks whose length is larger than the characteristic length of the correlations, and these blocks are then re-sampled to generate the bootstrap sample. In early experiments, we tried a form of ‘block bootstrap’ in which we filled each octant of a simulation cube with a block of data selected randomly from the cube. Unfortunately, the method did not work well. The sharp edges of the octants introduced spurious small-scale power, and the covariance of small-scale power differed substantially from that measured by the ensemble method. The mathematical relation between the covariance of power obtained by the block bootstrap method and the true covariance of power is sufficiently obscure that it was difficult to assess the possible causes of the discrepancy.

In HRS, we argue that all variations on jackknife and bootstrap are really just different ways of re-weighting the data to yield a new estimate of the quantity of interest, and that the best way to avoid unpleasant side-effects on spatial data is to re-weight with a smooth function of position. Further, by re-weighting the simulation with a function that contains only large-scale Fourier modes, unpleasant numerical artefacts in the power spectrum are confined to the largest scales, making them much easier to deal with.

4.1 Covariance of power of weighted density

Following HRS, we define the i ’th weighted density by

$$\rho_i(\mathbf{r}) \equiv w_i(\mathbf{r})\rho(\mathbf{r}). \quad (17)$$

We use the minimum variance set of weightings recommended in section 3 of HRS. In real space the i ’th weighting has the form

$$w_i(\mathbf{r}) = \sqrt{2} \cos \left[2\pi \left(\mathbf{k}_i \cdot \mathbf{r} + \frac{1}{16} \right) \right], \quad (18)$$

where

$$\mathbf{k}_i = \begin{cases} \{1, 0, 0\} & 12 \text{ weightings} \\ \{1, 1, 0\} & 24 \text{ weightings} \\ \{1, 1, 1\} & 16 \text{ weightings.} \end{cases} \quad (19)$$

The different weightings are obtained by all possible reflections and rotations of the components of \mathbf{k} , which yields 26 weightings in total, allowing for all of the symmetries in equation (18). A further 26 weightings are obtained by adding a phase shift of $\pi/2$, i.e. $1/16 \rightarrow 5/16$ in equation (18), equivalent to translating one of the coordinates by a quarter box (because of the symmetry of the weighting functions, only one such translation yields a distinct weighting). The estimate of power from this second set of weightings is predicted by HRS to be highly anti-correlated with that obtained from the first set of 26 weightings, but they are sufficiently uncorrelated that including all 52 weightings does yield a better measurement of the covariance matrix than is obtained from only 26 weightings. It is possible to apply more weightings constructed from higher order modes than those in equation (19) but, as argued in section 3.5 of HRS, these contain progressively less independent information, and yield progressively less accurate estimates of covariance.

A covariance matrix estimated as the average of n distinct estimates can have a rank no greater than n , a fact

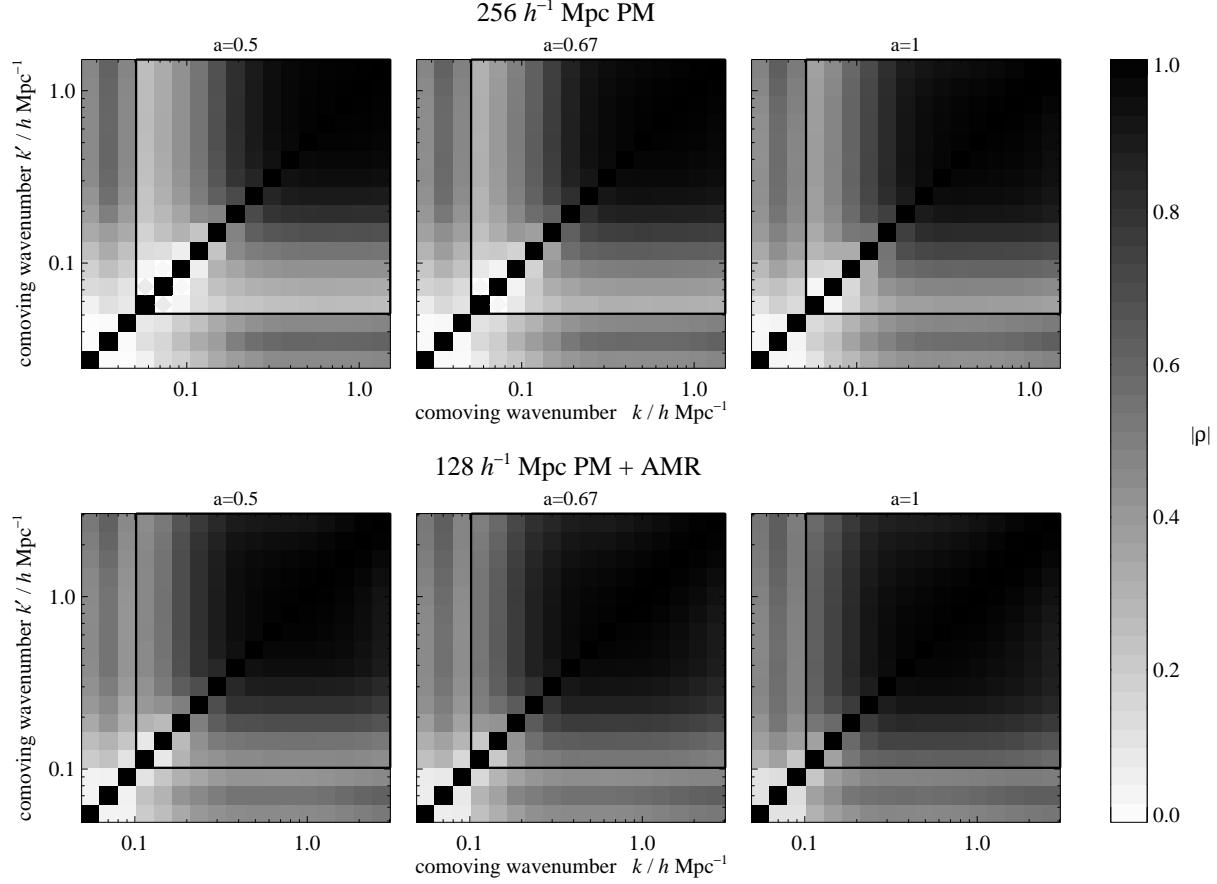


Figure 2. Correlation matrices of estimates of non-linear power using the re-weighting scheme of HRS at 3 epochs (left to right $a = 0.5$, 0.67 and 1). Correlation matrices were estimated separately for each simulation and then averaged over simulations. Results shown are averages over 100 PM simulations of each box size and additionally, in the case of the $128 h^{-1} \text{ Mpc}$ boxes, the 25 ART simulations. Greyscale is used to indicate the magnitude of the correlations, ranging from 0 (no correlation) to 1 (perfect correlation). A heavy, black border outlines the region of each covariance matrix that is unaffected by numerical artefacts from the re-weighting scheme. Bins outside of the bordered area are excluded from further analyses.

also pointed out recently by Pan & Szapudi (2005). The 52 weightings given by equation (19) prove sufficient to yield, for each simulation, a non-singular (no zero eigenvalues), and indeed positive definite (no negative eigenvalues) estimate of the covariance matrix for the 20 bins of wavenumber used here.

As a practical matter, we implement the re-weighting scheme by weighting individual particles, before assigning the density to the Fourier mesh. The weighted overdensity at point j on the mesh is defined to be

$$\delta_i(\mathbf{r}_j) \equiv \frac{\rho_i(\mathbf{r}_j)}{\bar{\rho}} - 1, \quad (20)$$

where $\bar{\rho}$ is the mean of the *unweighted* density field. Note that this is a different convention to that used in HRS, in which the quantity being transformed is

$$\Delta\rho_i(\mathbf{r}) \equiv \rho_i(\mathbf{r}) - \bar{\rho}_i(\mathbf{r}), \quad (21)$$

where $\bar{\rho}_i(\mathbf{r}) = w_i(\mathbf{r})$ and $\bar{\rho} \equiv 1$. The difference between these two conventions only affects the power on the largest scales (those wavebands containing modes appearing in equation 19) and can most easily (and accurately) be corrected for in Fourier space. However, since the covariance on

these scales is not correctly reproduced by the re-weighting method, we simply exclude them from our analyses.

The covariance of power over the ensemble of weighted powers is related to the true covariance of power by (HRS)

$$\langle \Delta\hat{P}(k)\Delta\hat{P}(k') \rangle = 2\langle \Delta\hat{P}_i(k)\Delta\hat{P}_i(k') \rangle_i, \quad (22)$$

where $\langle \dots \rangle_i$ denotes an average over different weightings and the factor 2 corrects the covariance of the ensemble to the true covariance of estimates of power. The deviation $\Delta\hat{P}_i(k)$ in the power spectrum of the i 'th weighted density must be measured relative to some expected or mean value, and HRS discussed two possibilities. The first is to measure the deviation relative to the power spectrum of the unweighted density of the simulation:

$$\Delta\hat{P}_i(k) \equiv \hat{P}_i(k) - \hat{P}(k). \quad (23)$$

The second is to measure the deviation relative to the mean of the power spectra of the weighted densities:

$$\Delta\hat{P}_i(k) \equiv \hat{P}_i(k) - \frac{1}{N} \sum_i \hat{P}_i(k). \quad (24)$$

The advantage of the first strategy, equation (23), is that the power spectrum of the unweighted density is, by symmetry,

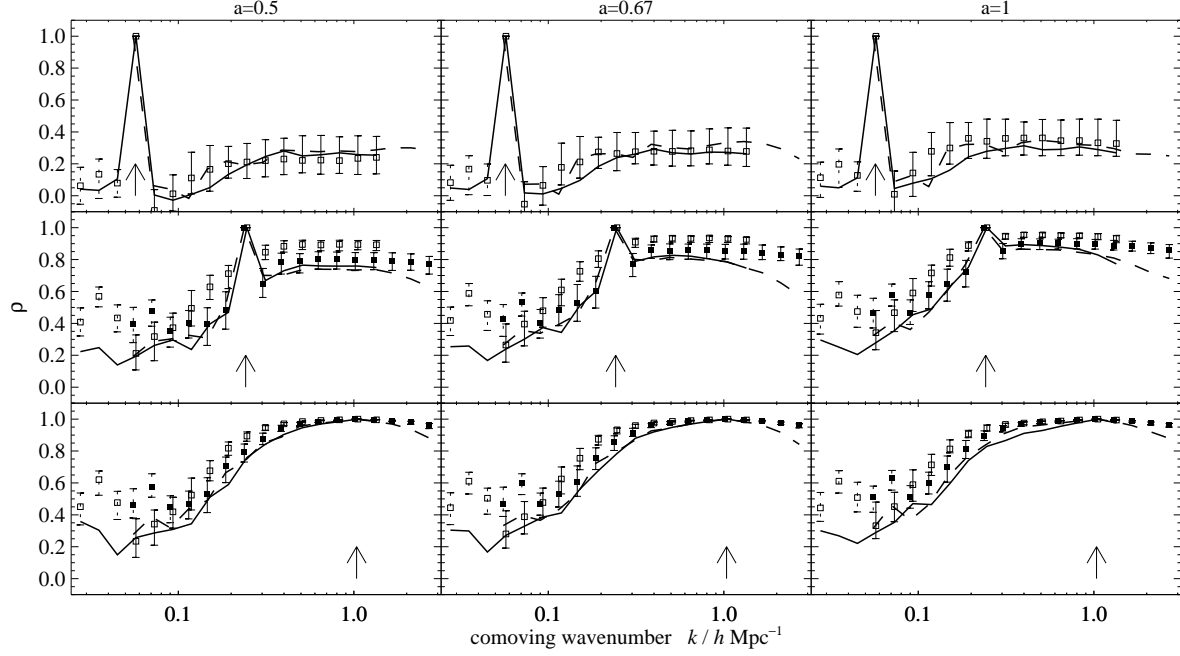


Figure 3. Correlations between estimates of the non-linear power spectrum using the re-weighting method, with ΔP_i defined by equation (24). The three columns show cross-sections through the correlation matrices at 3 epochs (left to right $a = 0.5, 0.67$ and 1), while each row shows a cross-section at a different wavenumber ($k' = 0.057, 0.242$ and $1.036 h \text{ Mpc}^{-1}$), marked by the vertical arrow in each panel. Symbols with error bars mark the median and quartiles of the distribution of correlation co-efficients measured from each individual simulation for the $256 h^{-1} \text{ Mpc}$ PM (open symbols) and $128 h^{-1} \text{ Mpc}$ PM+ART (filled symbols) simulations. Dotted error bars are used for points that are outside of the range of wavenumbers for which the covariance is expected to be reliably estimated by the re-weighting method. Solid and dashed lines show the correlations between (unweighted) estimates of power using the full ensemble of $256 h^{-1} \text{ Mpc}$ PM and $128 h^{-1} \text{ Mpc}$ PM+ART simulations, respectively. Data for the $128 h^{-1} \text{ Mpc}$ boxes are missing in the top row because this entire cross-section lies outside the reliable region for this box size.

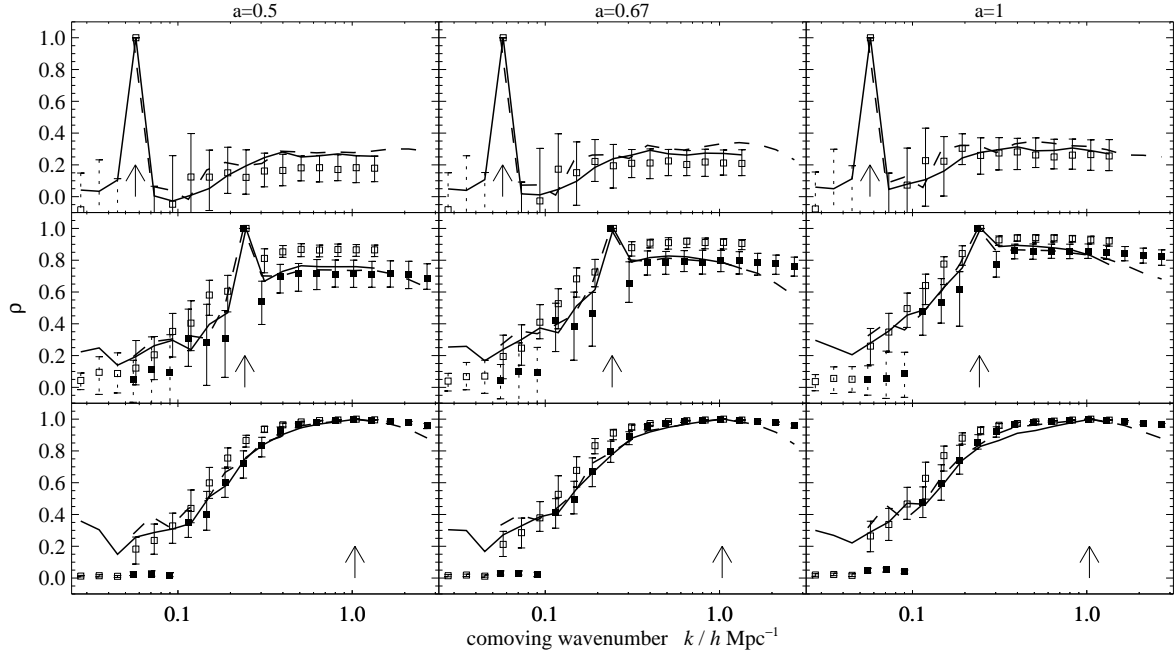


Figure 4. As Fig. 3, but with ΔP_i defined by equation (23). The symbols and lines have the same meanings as in Fig. 3. Together, the two figures demonstrate that the different methods give broadly consistent results.

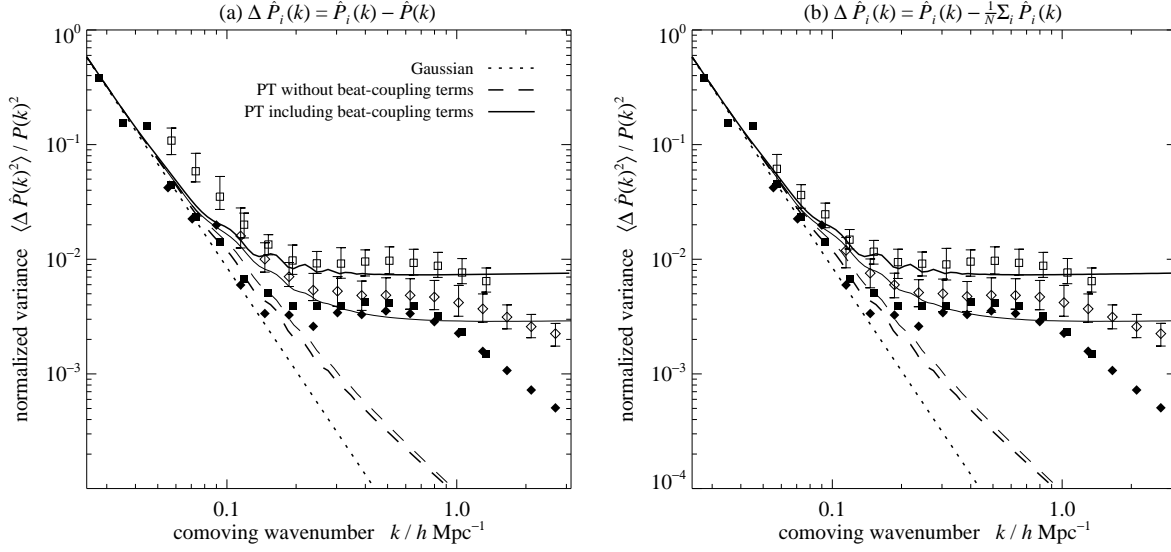


Figure 5. Variance of estimates of power using the re-weighting method with ΔP_i defined by (a) equation (24) and (b) equation (23). Open symbols show the variance, normalized to the square of the power, at $a = 1$ in the $256 h^{-1}$ Mpc simulations (squares) and the $128 h^{-1}$ Mpc simulations (diamonds). The results for the $128 h^{-1}$ Mpc box size have been shifted vertically by a factor 8 to compensate for the reduced density of modes and enable direct comparison between the two box sizes. Error bars mark the upper and lower quartiles of the distribution of estimates of variance over individual simulations. Solid symbols give the corresponding results measured using the ensemble method (these points are the same in both panels). The solid and dashed curves are predictions from perturbation theory, both with (solid) and without (dashed) the beat-coupling terms (see HRS for details of the calculation). The heavier lines are the results for the $256 h^{-1}$ Mpc box size. The dotted line is the expected variance for Gaussian fluctuations.

the most accurate estimate of power in a simulation, so the statistical uncertainty is potentially least in this case. However, the power spectra of weighted densities are likely to be slightly biased relative to the power spectrum of the unweighted density, because weighting the density effectively smooths the power, which biases it if power is other than a linear function of wavenumber. The advantage of the second strategy, equation (24), is that it removes this slight bias, so the systematic uncertainty is potentially smaller in this case.

The lower panel of Fig. 1 shows the median and quartiles of the distribution of the deviations between the averaged power spectra of weighted densities and the power spectrum of the unweighted density of each simulation (c.f. equation 20 of HRS). The two agree well on small scales, but on larger scales they can differ by up to 20 percent in extreme cases.

In the following sections, we show results for both strategies, equations (23) and (24), and find that the two strategies give consistent results. However, we do find that equation (23) gives variances that are slightly but systematically higher than those from equation (24), which we attribute to the systematic bias between the power spectra of weighted densities versus the power spectrum of the unweighted density. For the purposes of computing information, Section 5, we therefore choose the second strategy, equation (24).

Each set of weightings in equation (19) is generated from a different wavenumber $k_i = |\mathbf{k}_i|$, so we expect estimates of power from each set to be biased in a slightly different way. This bias is small; nevertheless it is preferable to estimate the covariance matrix separately from each set of weightings and then combine them to obtain a single es-

timate of the covariance matrix, weighting by the number of weightings in each set. This is the procedure that we adopt in Section 5.

4.2 Tests

In this section we describe several tests of the measurement of covariance of power. In Section 4.2.1 we show that the weightings and ensemble methods give consistent results for the correlation coefficients of band-powers. By contrast, in Section 4.2.2 we show that the two methods give substantially different results for the variance of power at non-linear scales. In Section 4.2.3 we check the assumption that the distribution of estimates of power is (thanks to the central limit theorem) adequately Gaussian.

4.2.1 Correlations in the power spectrum

Fig. 2 shows the matrix of correlation coefficients,

$$\rho_{kk'} \equiv \frac{\langle \Delta P_k \Delta P_{k'} \rangle}{\sqrt{\langle \Delta P_k^2 \rangle \langle \Delta P_{k'}^2 \rangle}}, \quad (25)$$

of estimates of the non-linear power spectrum at three epochs, measured using the re-weighting scheme outlined in Section 4.1. Each matrix is the average result from 100 individual PM simulations and, in the case of the $128 h^{-1}$ Mpc boxes, 25 ART simulations. The final epoch ($a = 1$) can be directly compared to fig. 2 of Paper I, in which we show the results from the ensemble method. We expect numerical artefacts from the re-weighting to be restricted to wavenumbers $k \leq \sqrt{3}k_b$, the highest wavenumber used in the weighting functions. Fig. 2 shows that this is indeed the case: the

degree of correlation changes abruptly between bins containing wavenumbers inside and outside this limit. In the following analysis, we use only bins with wavenumbers a factor of 2 away from the fundamental mode of the box, $k \geq 2k_b$.

Fig. 3 shows three cross-sections through each of the correlation matrices in Fig. 2. The data plotted are the medians and quartiles of the distribution over the independent realizations. The measurements from the re-weighting and ensemble methods are generally consistent, although the re-weightings method tends to yield somewhat higher correlations at smaller scales, and higher for the $256 h^{-1}\text{Mpc}$ boxes than the $128 h^{-1}\text{Mpc}$ boxes. The scatter between individual simulations is considerable, especially where the correlation coefficient is small.

Fig. 4 shows the same results, but with the deviations $\Delta\hat{P}_i(k)$ in the power spectra of weighted densities being measured relative to the power spectrum of the unweighted density, equation (23), as opposed to the mean of the power spectra of weighted densities, equation (24). The correlation coefficients are similar to those shown in Figure 3, as they should be.

4.2.2 Variance of the power spectrum

While the re-weighting and ensemble methods yield consistent results for the correlation matrix of non-linear power, the variance of (and hence the information contained in) the non-linear power spectrum is an altogether different matter.

Fig. 5 shows the variance, normalized to the square of the unweighted power spectrum, estimated using the re-weighting method. We show results for both equation (23) and equation (24). Notice that, as expected, equation (23) overestimates the variance of power on large scales. On small scales, however, the results are entirely consistent.

For Gaussian fluctuations, the normalized variance equals $2/N_k$ (equation 5), where N_k is the number of modes in a given wavenumber bin. For a periodic box, the number of modes in a set of logarithmically-spaced bins increases with central wavenumber as $N_k \propto k^3$. The $128 h^{-1}\text{Mpc}$ simulations have fewer modes (by a factor of 8) at a given wavenumber, so the results for this box size have been shifted vertically down by this factor to allow a direct comparison between the results for the two box sizes. As in Fig. 3, the data shown in Fig. 5 are medians over many individual simulations, with error bars marking the quartiles of the distribution. The results for the $128 h^{-1}\text{Mpc}$ PM simulations and the $128 h^{-1}\text{Mpc}$ ART simulations are consistent, so we combine them into a single set of points for clarity (a figure showing the ART results alone appears as fig. 2 of HRS).

The two different box sizes yield consistent results when the ensemble method is used. For the re-weighting method, on the other hand, there are clear discrepancies, both between the re-weighting method and the ensemble method and between the two box sizes. At translinear and non-linear scales, the variance measured by re-weighting is significantly higher than that measured for the ensemble, particularly for the larger box size, and the discrepancy grows ever larger at smaller scales.

Fig. 5 also shows the predictions of perturbation theory. The calculation of these curves is described in section 4.3 of HRS. Perturbation theory helps to explain the discrepancies both between the results of the ensemble and re-weighting

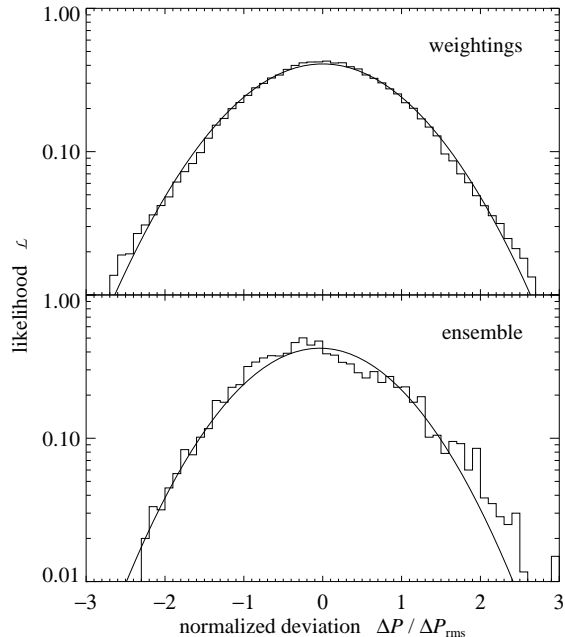


Figure 6. Distributions of estimates of non-linear power, equation (24), using the re-weighting method (top panel) and the ensemble method (bottom panel). Individual histograms for all bands with wavenumbers $k > 0.2 h^{-1}\text{Mpc}$, containing at least 2000 Fourier modes, are scaled to unit variance and stacked. The smooth curves are Gaussian fits to the data, assuming Poisson weighting of the counts in each bin.

methods, and between the two different box sizes. The variance measured by the re-weighting technique departs from the ensemble result where the (constant) beat-coupling term (equation 94 of HRS) becomes the dominant source of variance. The source of this term – coupling of closely-spaced Fourier modes to the large-scale beat mode between them – is discussed in section 4 of HRS. The magnitude of the beat-coupling term is proportional to the amplitude of the power spectrum on roughly the size of the box, which explains why different sizes of simulation box yield systematically different estimates of the small-scale variance. Perturbation theory correctly predicts the magnitude of the discrepancy between the small-scale variance estimated from the two different sizes of simulation.

Note that perturbation theory fails to reproduce the correct non-linear variance for the ensemble method (dashed lines), presumably because perturbation theory breaks down in the highly non-linear regime. This may be responsible for the small discrepancies between the results of the re-weightings method and the theoretical curves in Fig. 5.

4.2.3 Distribution of estimates of power

In Section 2 we asserted that, thanks to the central limit theorem, estimates of power should be Gaussianly distributed about their expectation value, even in the highly non-linear regime. Since our analysis relies on the validity of this assumption it is worth putting to the test.

Fig. 6 shows the distribution of deviations of estimates of power using both the re-weighting method and the en-

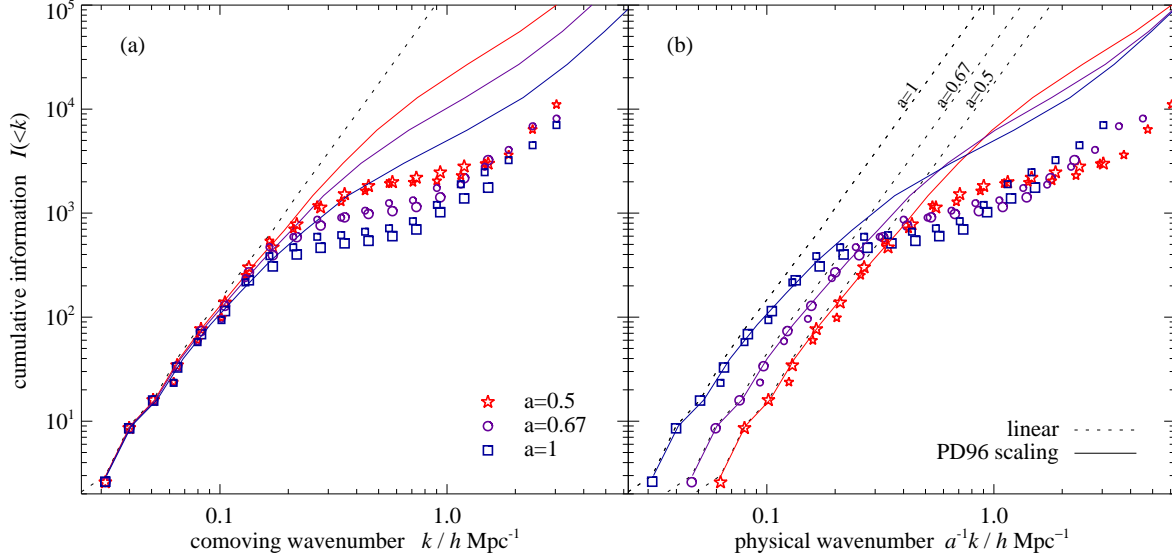


Figure 7. Cumulative information in the non-linear power spectrum at 3 epochs (top to bottom: $a = 0.5, 0.67$ and 1), as a function of (a) comoving and (b) physical wavenumber. Large symbols are points derived from the $256 h^{-1}$ Mpc PM simulations and small simulations are for the $128 h^{-1}$ Mpc PM+ART boxes. The results for the $128 h^{-1}$ Mpc boxes have been shifted vertically by a factor 8 to account for the higher density of modes at a given comoving k and allow direct comparison with the larger box size. The dotted line marks the expected amount of information for Gaussian fluctuations and the solid curves are the result of applying the PD96 scaling to the dotted line at each epoch. This is a revised version of fig. 3 in Paper I (see text for details).

semble method for wavenumbers in the non-linear regime (which we take to be $k > 0.2 h^{-1}$ Mpc). Individual histograms for each waveband have been scaled to unit variance and stacked to produce a single distribution for each method. The estimates of power from the re-weighted simulations are indeed distributed close to Gaussianly. Assuming Poisson statistics, the value of χ^2 for the fit is 5.84 per degree of freedom, which seems reasonable, given the high degree of correlation between estimates of non-linear power. For the ensemble method, the distribution is also close to Gaussian, although there are deviations from Gaussianity – in particular the presence of a sharper than Gaussian peak and a tail of large, positive deviations that cause the actual mean and variance of power to be slightly larger than the fitted values. The value of χ^2 in this case is 3.86 per degree of freedom, for Poisson statistics.

5 INFORMATION CONTENT OF THE NON-LINEAR POWER SPECTRUM

In the previous section, we showed that measuring the covariance of non-linear power by re-weighting an individual simulation yields substantially different results at non-linear scales than is found from the scatter over an ensemble of simulations. The discrepancy is consistent with the explanation proposed by HRS: beat-coupling between the covariance on non-linear scales and the power on large scales. The difference in covariance between the two methods translates directly into a difference in the quantity of information in the non-linear power spectrum.

5.1 Method

We decorrelate estimates of the (log-) amplitude for each simulation individually, using the covariance matrix estimated using the re-weighting method. As our fiducial power spectrum P_k in equation (9) we use the true (unweighted) power in each simulation. The re-weighting method restricts the range of wavenumbers for which the covariance matrix – and hence the Fisher information – can be reliably measured to $k > \sqrt{3}k_b$. Since our purpose is to measure the *cumulative* information, equation (7), up to some wavenumber k_{\max} , the contribution from wavenumbers smaller than this limit must be taken into account. For the $256 h^{-1}$ Mpc boxes we assume that the excluded bins contain the expected amount of information for Gaussian fluctuations ($N_k/2$, where N_k is the number of Fourier modes in those bins). That this is a reasonable assumption is confirmed by the fact that the first few bins for which we do have measurements of the variance using the re-weighting method follow the Gaussian expectation closely (see the top panel of Fig. 5). For the $128 h^{-1}$ Mpc boxes there is a further complication. The lower limit on the wavenumbers accessible using the re-weighting method brings us into the regime in which non-linear effects start to become important. The most reasonable thing to do here would seem to be to use the results from the $256 h^{-1}$ Mpc boxes to estimate the quantity of missing information. Although there are clearly systematic differences between the results of the re-weighting method for different box sizes, these are small at the scales in question.

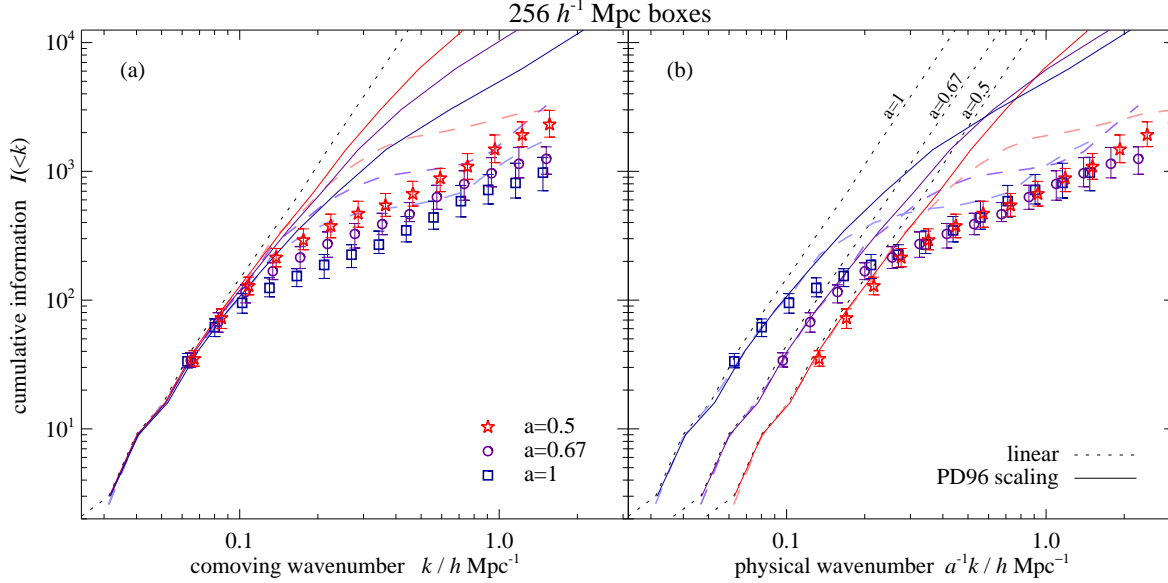


Figure 8. Cumulative information, as a function of (a) comoving and (b) physical wavenumber, for the same 3 epochs as Fig. 7. Points with error bars are medians and quartiles of the results measured using the re-weighting method on 100 PM simulations with a box size of $256 h^{-1}$ Mpc. For clarity, the three sets of points have been artificially separated by a small factor in wavenumber, with the $a = 0.67$ points having the correct wavenumber. For comparison, the ensemble results from Fig. 7 are shown as light, dashed lines. The dotted line marks the expected amount of information for Gaussian fluctuations and the solid curves are the result of applying the PD96 scaling to the dotted line at each epoch.

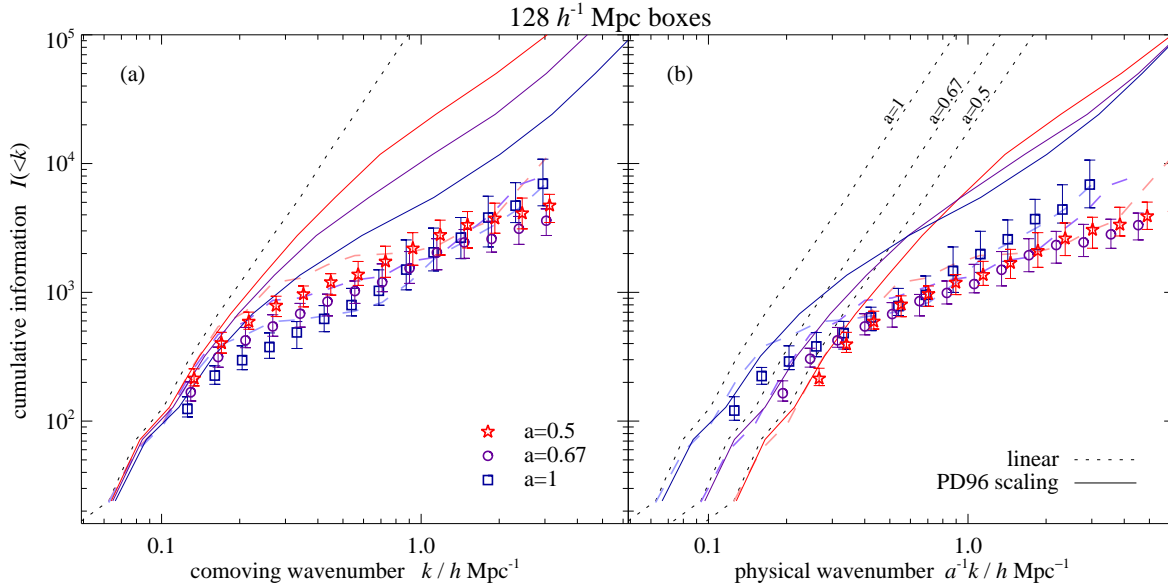


Figure 9. Cumulative information, as Fig. 8, for 100 PM simulations and 25 ART simulations with a box size of $128 h^{-1}$ Mpc. The results have been shifted vertically by a factor of 8 to enable direct comparison with the larger box size. The symbols and lines have the same meanings as in Fig. 8, but note the different axis ranges.

5.2 Results

5.2.1 Ensemble method

Fig. 7 shows the cumulative information as a function of wavenumber at 3 epochs ($a = 0.5, 0.67$ and 1) for the ensemble method of estimating the covariance of power.

The curves in Fig. 7 differ somewhat from those in fig. 3

of Paper I. In the previous paper, we (incorrectly) decorrelated the power spectrum *prior* to multiplying by the partial derivatives in equation (7). As discussed in Section 2.4, this is only a good approximation to the exact result provided that the band-power windows are narrow in k , which is a poor assumption in our case because of the existence of broad correlations in the power spectrum. The curves in

Fig. 7 (and Figs. 8 and 9 later) were produced using the exact formalism set out in Section 2.4 of the present paper. We still find that the ensemble method yields very little independent information in the translinear regime ($k \simeq 0.2\text{--}0.8\ h^{-1}\text{ Mpc}$), over and above that in the linear regime, although the flatness is not as pronounced as suggested by our previous results. In fact, the cumulative information increases by a factor of approximately 2 over the aforementioned range of wavenumbers (at $a = 1$), still much less than the factor of $k^3 \simeq 64$ expected for linear fluctuations. There remains a sudden upturn in the cumulative information on small scales, implying that the power spectrum at fully non-linear scales contains unique information about the amplitude of the initial power spectrum, that is not found in the present day linear power spectrum.

In Paper I, we interpreted the decrease in information in the translinear regime as the result of rapid transfer of information from larger to smaller scales. An alternative explanation, which was mentioned in Paper I but discarded as being contrived, is that information is temporarily diverted into higher order statistics, such as the bispectrum, in the translinear regime. Since filamentary structures are more readily described by higher-order statistics than by the power spectrum alone, it is entirely plausible that, on translinear scales, the bispectrum does contain information not present in the power spectrum, about the initial conditions of structure formation, and that this information is somehow returned to the power spectrum on smaller scales.

We argued in Paper I that, if information is conserved overall then, under the assumption of stable clustering, the amount of information up to a given physical (as opposed to comoving) wavenumber ought to be independent of time in the fully non-linear regime. The right panel of Fig. 7 shows the cumulative information for the same three epochs, plotted as a function of physical wavenumber k/a . The results are consistent with the hypothesis that information is largely conserved by non-linear evolution.

5.2.2 Re-weighting method

In Figs. 8 and 9 we compare the results from the ensemble method with those from the re-weighting method. For the re-weighting method we show the median and quartiles of the distribution of results from the individual simulations. Qualitatively, the behaviour of the cumulative information – as a function of both wavenumber and cosmic epoch – for the re-weighting method is similar to that for the ensemble method. However, the flattening in the translinear regime is less pronounced than for the ensemble case and there is no clear evidence for an upturn on small scales, although the curves for the re-weighting method follow the average slope of those from the ensemble method. Overall, the information measured using the re-weighting method is considerably less than when the ensemble method is used. Beat-coupling to large scales prevents much of the information that is, in principle, contained in the power spectrum from being extracted when the re-weighting method is used. It is also worth noting that the magnitude of the beat-coupling effect depends on the size of the simulation, so that the two different box sizes are no longer in complete agreement on small scales. The $128\ h^{-1}\text{ Mpc}$ boxes (Fig. 9) are in closer agreement with the ensemble method, as expected.

As predicted by perturbation theory, the beat-coupling contribution to the covariance, which is a factor $\sim P(2k_b)/P(k)$ larger than the other terms at a given k , becomes increasingly dominant at smaller scales. For the power spectrum considered here, the contribution from beat-coupling also increases with cosmic epoch. We would therefore not expect the cumulative information, measured using the re-weighting method, up to a given physical wavenumber to be necessarily constant over time, even in the stable clustering regime.

6 SUMMARY

This paper extends and expands on the results reported in Paper I concerning the Fisher information contained in the non-linear power spectrum about the amplitude of the initial (post-recombination) linear power.

In Paper I, we measured the covariance of power from the scatter in power over a large ensemble of simulations. Here we reported measurements of covariance of power from both the ensemble method and a new method, described in a companion paper (HRS), in which smoothly varying weighting functions are applied to each simulation to yield a separate estimate of the covariance of power for each simulation.

We have shown that the two methods yield substantially different estimates for the covariance of power at non-linear scales. This does not mean, however, that one or other of the methods is incorrect. Rather, it turns out that measuring the covariance of power is a more subtle problem than we had previously suspected. Beat-coupling – the process whereby the covariance between Fourier modes separated by a small wavevector couple by gravitational growth to the large-scale beat mode between them – dominates the covariance on non-linear scales. We compared our results to a calculation using perturbation theory (HRS) and found that the theory explains qualitatively the features of beat-coupling, seen in the simulations.

Beat-coupling contributions to the covariance of power occur whenever Fourier modes have a finite width, as opposed to being delta-functions at discrete wavevectors. As argued by HRS, this means that beat-coupling is likely to be important in real galaxy surveys. The ensemble method, on the other hand, measures covariances between the amplitudes of Fourier modes with precisely defined (delta function) wavenumbers. If information is conserved by non-linear evolution then it is this quantity that, overall, remains invariant with time.

Theory predicts that the effects of beat-coupling will be greatest when the largest scales in a survey are close to the peak of the power spectrum ($k \simeq 0.016\ h\text{ Mpc}^{-1}$; $\pi/k \sim 200\ h^{-1}\text{ Mpc}$). If this is true then our results suggest the covariance of power in such a survey will be dominated by beat-coupling on small scales and, counter-intuitively, will be sensitive to the power on the largest scales in the survey, leading to a reduction in the amount of information extractable from the power spectrum. As pointed out in HRS, the best way to test this is using mock galaxy catalogues drawn from a single large simulation using the same selection function as the survey observations.

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